

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100423\
 Data File : BM042218.D
 Acq On : 05 Oct 2023 15:53
 Operator : MA/JU
 Sample : PB155938BS
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS938

Quant Time: Oct 06 05:27:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 06 05:25:26 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.929	152	3537	0.400	ng/ul	0.00
4) Naphthalene-d8	10.741	136	9519	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.578	164	3984	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.329	188	9105	0.400	ng/ul	#-0.01
17) Chrysene-d12	21.507	240	4595	0.400	ng/ul	0.00
23) Perylene-d12	23.909	264	4648	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.356	96	3466	0.736	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.330	152	4288	0.328	ng/ul	0.00
18) Fluoranthene-d10	19.352	212	7396	0.429	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.393	88	8239	1.675	ng/ul#	77
5) Naphthalene	10.790	128	9782	0.308	ng/ul	99
7) 2-Methylnaphthalene	12.401	142	5524	0.291	ng/ul	95
8) 1-Methylnaphthalene	12.621	142	5763	0.301	ng/ul	97
10) Acenaphthylene	14.305	152	7901	0.314	ng/ul	99
11) Acenaphthene	14.643	153	6423	0.337	ng/ul	99
12) Fluorene	15.628	166	7359	0.343	ng/ul	96
14) Pentachlorophenol	16.966	266	2338	0.566	ng/ul	99
15) Phenanthrene	17.371	178	12502	0.333	ng/ul	99
16) Anthracene	17.464	178	10484	0.305	ng/ul	99
19) Fluoranthene	19.385	202	14298	0.410	ng/ul	97
20) Pyrene	19.752	202	15551	0.419	ng/ul	97
21) Benzo(a)anthracene	21.489	228	10980	0.370	ng/ul	99
22) Chrysene	21.545	228	11745	0.407	ng/ul	100
24) Benzo(b)fluoranthene	23.167	252	12203	0.445	ng/ul	98
25) Benzo(k)fluoranthene	23.217	252	11791	0.437	ng/ul	98
26) Benzo(a)pyrene	23.798	252	9997	0.423	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.398	276	13938	0.416	ng/ul#	88
28) Dibenzo(a,h)anthracene	26.424	278	10287	0.407	ng/ul	94
29) Benzo(g,h,i)perylene	27.169	276	12476	0.446	ng/ul	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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